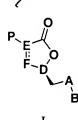
Claims

What is claimed is:

5 1. A compound of formula I:



or a pharmaceutically acceptable salt thereof, wherein:

10 A is O,

NH, or

S;

B is

15 $C(=O)R_1$,

 $C(=S)R_1$,

heterocylco,

heteroaryl,

C(=O)-heterocyclo,

C(=N)-CN, or

C(=O)-heteteroaryl;

either D is N, E is C, and F is CH when "-----" is a bond, or D is CH, E is N, and F is CH_2 when "-----" is absent;

P is

25

11 1

attachment;

5

J, K, Q independently are CR_2 or N, with the proviso that when any one of J, K, or Q is N, then the other two are CR_2 ;

"-----"is absent; or is a bond; and

X, Y, Z independently are C=C-R₅,

O=C,

10

CH₂,

 CHR_{3}

CHR₄,

CR₃R₄,

NR₅,

15

 $N(C=O)R_5$,

 $N(C=O)OR_5$,

NSO₂R₅,

NSO₂NR₅,

Ο,

20

S,

SO, or

SO₂;

 R_1 is H,

25

(C₁-C₈)alkyl,

(C₃-C₆)cycloalkyl,

O— $(C_1$ - C_4)alkyl,

O-(C₃-C₆)cycloalkyl,

 $S-(C_1-C_4)$ alkyl, S—(C₃-C₆)cycloalkyl, NH_2 5 $NH(C_1-C_4)$ alkyl, $N((C_1-C_4)alkyl)_2$, or NH—(C₃-C₆)cycloalkyl; R₂ is H, 10 halo, (C₁-C₈)alkyl, (C₃-C₆)cycloalkyl, O— $(C_1$ - C_4)alkyl, O—(C₃-C₆)cycloalkyl, $S-(C_1-C_4)$ alkyl, 15 S— $(C_3$ - $C_6)$ cycloalkyl, NH_2 , $NH(C_1-C_4)alkyl$, $N((C_1-C_4)alkyl)_2$, or 20 NH—(C₃-C₆)cycloalkyl; R₃ and R₄ independently are halo, (C_1-C_8) alkyl, (C₃-C₆)cycloalkyl, 25 O— $(C_1$ - C_4)alkyl, O—(C₃-C₆)cycloalkyl, $S-(C_1-C_4)$ alkyl, S—(C₃-C₆)cycloalkyl, NH_2 30 $NH(C_1-C_4)alkyl$, $N((C_1-C_4)alkyl)_2$ NH—(C₃-C₆)cycloalkyl;

aryl,
$$(CH_2)_n\text{-aryl},$$
 heterocyclo,
$$(CH_2)_n\text{-heterocyclo},$$

$$(CH_2)_n\text{-heteroaryl, or}$$

$$(CH_2)_n\text{-heteroaryl,}$$
 wherein n is 0, 1, 2, or 3;

 R_5 is H, $(C_1\text{-}C_8)\text{alkyl},$ $(C_3\text{-}C_6)\text{cycloalkyl},$ aryl, $(CH_2)_n\text{-aryl},$ heterocyclo, $(CH_2)_n\text{-heterocyclo},$ heteroaryl, or $(CH_2)_n\text{-heteroaryl},$

20 2. The compound of claim 1 as designated in formula IA.

wherein n is as defined above.

3. The compound of claim 1 as designated in formula IB.

25

4. The compound of claim 1 as designated in formula IC.

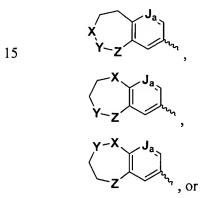
5

5. The compound of claim 1, wherein P is

10

wherein " $\mbox{\ensuremath{$\sim$}}\mbox{\ens$

6. The compound of claim 1, wherein P is



7. The compound of claim 6, wherein two of X, Y, or Z is $C=C-R_5$,

O=C,

5 NR₅,

 $N(C=O)R_5$,

 $N(C=O)OR_5$,

NSO₂R₅,

NSO₂NR₅,

10 O,

S,

SO, or

SO₂NR₅,

and the other of X, Y, or Z is CH₂ or CR₃R₄.

15

8. The compound of claim 7, wherein P is

20

9. The compound of claim 7, wherein P is

10. The compound of claim 7, wherein P is

11. The compound of claim 6, wherein one of X, Y, or Z is $C=C-R_5$,

O=C,

10 NR₅,

 $N(C=O)R_5$

 $N(C=O)OR_5$

NSO₂R₅,

NSO₂NR₅,

15 O,

S,

SO, or

SO₂NR₅,

and the other of X, Y, or Z is CH₂.

20

5

12. The compound of claim 11, wherein P is

5 13. A compound of formula II

or a pharmaceutically acceptable salt thereof, wherein:

B is
$$C(=O)R_1,$$

$$C(=S)R_1,$$
 heterocylco, heteroaryl,
$$C(=O)\text{-heterocyclo,}$$

$$20 \hspace{1cm} C(=N)\text{-CN,or}$$

$$C(=O)\text{-heteroaryl;}$$

either D is N, E is C, and F is CH when "-----" is a bond, or D is CH, E is N, and F is CH_2 when "-----" is absent;

```
J, K, Q independently are CR2 or N, with the proviso that when any
                    one of J, K, or Q is N, then the other two are CR<sub>2</sub>;
                               "-----"is absent; or is a bond; and
 5
                               X, Y, Z independently are C=C-R<sub>5</sub>,
                                          O=C,
                                          CH<sub>2</sub>,
                                          CHR<sub>3,</sub>
                                          CHR<sub>4</sub>,
10
                                          CR<sub>3</sub>R<sub>4</sub>,
                                          NR<sub>5</sub>,
                                          N(C=O)R_5
                                          N(C=O)OR_5,
                                          NSO<sub>2</sub>R<sub>5</sub>,
15
                                          NSO<sub>2</sub>NR<sub>5</sub>,
                                          Ο,
                                          S,
                                          SO, or
                                          SO<sub>2</sub>;
20
                               R<sub>1</sub> is H,
                                          (C_1-C_8)alkyl,
                                          (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
                                          O—(C_1-C_4)alkyl,
                                          O—(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
25
                                          S—(C_1-C_4) alkyl,
                                          S—(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
                                          NH<sub>2</sub>,
                                          NH(C<sub>1</sub>-C<sub>4</sub>)alkyl,
30
                                          N((C_1-C_4)alkyl)_2, or
```

NH—(C₃-C₆)cycloalkyl,

```
R<sub>2</sub> is H,
                                             halo,
                                             (C_1-C_8)alkyl,
                                             (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
  5
                                             O—(C_1-C_4)alkyl,
                                             O—(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
                                             S-(C_1-C_4) alkyl,
                                             S—(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
                                             NH<sub>2</sub>,
10
                                             NH(C_1-C_4)alkyl,
                                             N((C_1-C_4)alkyl)_2, or
                                             NH—(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl;
                                 R<sub>3</sub> and R<sub>4</sub> independently are halo,
15
                                             (C_1-C_8)alkyl,
                                             (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
                                             O—(C_1-C_4)alkyl,
                                             O—(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
                                             S—(C_1-C_4) alkyl,
20
                                             S—(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
                                             NH<sub>2</sub>,
                                             NH(C<sub>1</sub>-C<sub>4</sub>)alkyl,
                                             N((C_1-C_4)alkyl)_2,
                                             NH—(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl;
25
                                             aryl,
                                             (CH<sub>2</sub>)<sub>n</sub>-aryl,
                                             heterocyclo,
                                             (CH<sub>2</sub>)<sub>n</sub>-heterocyclo,
                                             heteroaryl, or
30
                                             (CH<sub>2</sub>)<sub>n</sub>-heteroaryl,
                                 wherein n is 0, 1, 2, or 3;
```

(C₁-C₈)alkyl, (C₃-C₆)cycloalkyl, aryl, (CH₂)_n-aryl, heterocyclo, (CH₂)_n-heterocyclo, heteroaryl, or

wherein n is as defined above.

13. The compound of claim 12 as designated in formula IIA.

(CH₂)_n-heteroaryl,

15

5

14. The compound of claim 12 as designated in formula IIB.

20 15. The compound of claim 12 as designated in formula IIC.

16. The compound of claim 12 as designated in formula IID.

wherein J_a is N or CR_6 , wherein R_6 is H or F.

10 17. The compound of claim 12 as designated in formula IIE.

18. The compound of claim 17, wherein two of X, Y, or Z is C=C-R₅,

O=C,

15 NR₅,

5

 $N(C=O)R_5$

 $N(C=O)OR_5$

NSO₂R₅,

 $NSO_{2}NR_{5}, \\$

20 O,

S,

.

SO, or

SO₂NR₅,

and the other of X, Y, or Z is CH2 or CR3R4.

5 19. The compound of claim 18 as designated in formula IIF.

20. The compound of claim 18 as designated in formula IIG.

10

21. The compound of claim 20 as designated in formula IIH.

15

IIH

22. The compound of claim 20, wherein one of X, Y, or Z is C=C-R₅,

O=C,

NR₅,

 $N(C=O)R_5$

 $N(C=O)OR_5$,

NSO₂R₅,

NSO₂NR₅,

Ο,

S,

5 SO, or

SO₂NR₅,

and the others of X, Y, or Z is CH₂.

23. A compound of formula III /

III

10

or a pharmaceutically acceptable salt thereof, wherein:

A is O,

NH, or

S;

B is $C(=O)R_1$,

 $C(=S)R_1$,

20 heterocylco,

heteroaryl,

C(=O)-heterocyclo,

C(=N)-CN,or

C(=O)-heteteroaryl;

25

either D is N, E is C, and F is CH when "-----" is a bond, or D is CH, E is N, and F is CH_2 when "-----" is absent;

J, K, Q independently are CR_2 or N, with the proviso that when any one of J, K, or Q is N, then the other two are CR_2 ;

```
"-----"is absent or is a bond;
  5
                                 X, Y, Z independently are C=C-R<sub>5</sub>,
                                             O=C,
                                             CHR<sub>3</sub>
                                             CHR<sub>4</sub>,
10
                                             CR<sub>3</sub>R<sub>4</sub>,
                                            NR<sub>5</sub>,
                                            N(C=O)R_5
                                             N(C=O)OR_5,
                                            NSO<sub>2</sub>R<sub>5</sub>,
15
                                            NSO<sub>2</sub>NR<sub>5</sub>,
                                             Ο,
                                            S,
                                             SO, or
                                            SO<sub>2</sub>;
20
                                 R_1 is H,
                                            (C_1-C_8)alkyl,
                                            (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
                                            O—(C_1-C_4)alkyl,
                                            O—(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
25
                                            S—(C_1-C_4) alkyl,
                                            S—(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
                                            NH<sub>2</sub>,
                                            NH(C<sub>1</sub>-C<sub>4</sub>)alkyl,
30
                                            N((C_1-C_4)alkyl)_2, or
                                            NH—(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl;
```

```
R<sub>2</sub> is H,
                                           halo,
                                           (C_1-C_8)alkyl,
                                          (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
                                          O—(C_1-C_4)alkyl,
  5
                                           O—(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
                                           S—(C_1-C_4) alkyl,
                                           S—(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
                                           NH<sub>2</sub>,
                                          NH(C_1-C_4)alkyl,
10
                                          N((C_1-C_4)alkyl)_2, or
                                          NH—(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl;
                               R<sub>3</sub> and R<sub>4</sub> independently are H,
15
                                          halo,
                                           (C_1-C_8)alkyl,
                                          (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
                                          O—(C_1-C_4)alkyl,
                                          O—(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
20
                                          S—(C_1-C_4) alkyl,
                                          S—(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
                                          NH<sub>2</sub>,
                                          NH(C_1-C_4)alkyl,
                                          N((C_1-C_4)alkyl)_2,
25
                                          NH—(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl;
                                          aryl,
                                          (CH_2)_n-aryl,
                                          heterocyclo,
                                          (CH<sub>2</sub>)<sub>n</sub>-heterocyclo,
30
                                          heteroaryl, or
                                          (CH<sub>2</sub>)<sub>n</sub>-heteroaryl,
                               wherein n is 0, 1, 2, or 3;
```

$$R_5$$
 is H,
$$(C_1\text{-}C_8)\text{alkyl},$$

$$(C_3\text{-}C_6)\text{cycloalkyl},$$

$$\text{aryl},$$

$$(CH_2)_n\text{-aryl},$$

$$\text{heterocyclo},$$

$$(CH_2)_n\text{-heterocyclo},$$

$$\text{heteroaryl}, \text{ or }$$

$$(CH_2)_n\text{-heteroaryl},$$

$$\text{wherein n is as defined above}.$$

24. The compound of claim 23 as designated in formula IIIA.

IIIA

25. The compound of claim 23 as designated in formula IIIB.

20

15

26. The compound of claim 23 as designated in formula IIIC.

27. The compound of claim 27 as designated in formula IIID.

5

wherein Ja is N or CR6, wherein R6 is H or F.

10 28. The compound of claim 27 as designated in formula IIIE.

29. The compound of claim 28, wherein two of X, Y, or Z is $C=C-R_5$,

15
$$O=C$$
, NR_5 , $N(C=O)R_5$, $N(C=O)OR_5$, NSO_2R_5 , NSO_2NR_5 ,

O,

S,

SO, or

SO₂NR₅,

and the other of X, Y, or Z is CH_2 or CR_3R_4 .

30. The compound of claim 29 as designated in formula IIIG.

IIIG

10

31. The compound of claim 29 as designated in formula IIIH.

IIIH

15 32. The compound of claim 28, wherein one of X, Y, or Z is $C=C-R_5$,

O=C,

NR₅,

 $N(C=O)R_5$

 $N(C=O)OR_5$,

NSO₂ R_5 ,

NSO₂NR₅,

O,

S,

SO, or

SO_2NR_5 , and the other of X, Y, or Z is CH_2 .

33. A compound of formula IV:

.

or a pharmaceutically acceptable salt thereof, wherein:

A is O,

NH, or

5

S;

B is

 $C(=O)R_1$

15 $C(=S)R_1$,

heterocylco,

heteroaryl,

C(=O)-heterocyclo,

C(=N)-CN,or

20 C(=O)-heteteroaryl;

either D is N, E is C, and F is CH when "-----" is a bond, or D is CH, E is N, and F is CH_2 when "-----" is absent;

J, K, Q independently are CR₂ or N, with the proviso that when any one of J, K, or Q is N, then the other two are CR₂;

"-----"is absent; or is a bond; and

X, Y, Z independently are C=C-R₅,

```
O=C,
                                               CH<sub>2</sub>,
                                               CHR<sub>3,</sub>
                                               CHR<sub>4</sub>,
  5
                                               CR<sub>3</sub>R<sub>4</sub>,
                                               NR<sub>5</sub>,
                                               N(C=O)R_5
                                               N(C=O)OR<sub>5</sub>,
                                               NSO<sub>2</sub>R<sub>5</sub>,
10
                                               NSO<sub>2</sub>NR<sub>5</sub>,
                                               O,
                                               S,
                                               SO, or
                                               SO<sub>2</sub>;
15
                                   R<sub>1</sub> is H,
                                               (C_1-C_8)alkyl,
                                               (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
                                               O-(C_1-C_4)alkyl,
20
                                               O—(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
                                               S—(C_1-C_4) alkyl,
                                               S—(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
                                               NH<sub>2</sub>,
                                               NH(C_1-C_4)alkyl,
25
                                               N((C_1-C_4)alkyl)_2, or
                                               NH—(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
                                  R<sub>2</sub> is H,
                                               halo,
30
                                               (C_1-C_8)alkyl,
                                               (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
                                               O—(C_1-C_4)alkyl,
```

O—(C₃-C₆)cycloalkyl,

```
S-(C_1-C_4) alkyl,
                                           S—(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
                                           NH_2,
  5
                                           NH(C_1-C_4)alkyl,
                                           N((C_1-C_4)alkyl)_2, or
                                           NH—(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl;
                                R<sub>3</sub> and R<sub>4</sub> independently are halo,
10
                                           (C_1-C_8)alkyl,
                                           (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
                                           O—(C_1-C_4)alkyl,
                                           O—(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
                                           S—(C_1-C_4) alkyl,
                                           S—(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
15
                                           NH<sub>2</sub>,
                                           NH(C_1-C_4)alkyl,
                                           N((C_1-C_4)alkyl)_2,
                                           NH—(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl;
20
                                           aryl,
                                           (CH_2)_n-aryl,
                                           heterocyclo,
                                           (CH<sub>2</sub>)<sub>n</sub>-heterocyclo,
                                           heteroaryl, or
                                          (CH<sub>2</sub>)<sub>n</sub>-heteroaryl,
25
                               wherein n is 0, 1, 2, or 3;
                               R<sub>5</sub> is H,
                                           (C_1-C_8)alkyl,
30
                                          (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
                                           aryl,
                                           (CH<sub>2</sub>)<sub>n</sub>-aryl,
```

heterocyclo, $(CH_2)_n$ -heterocyclo, heteroaryl, or $(CH_2)_n$ -heteroaryl,

5 wherein n is as defined above.

34. The compound of claim 33 as designated in formula IVA.

10

35. The compound of claim 33 as designated in formula IVB.

15 36. The compound of claim 33 as designated in formula IVC.

37. The compound of claim 33 as designated in formula IVD.

wherein J_a is N or CR_6 , wherein R_6 is H or F.

5

38. The compound of claim 33 as designated in formula IVE.

10 39. The compound of claim 38, wherein two of X, Y, or Z is $C=C-R_5$,

O=C,

NR₅,

 $N(C=O)R_5$

 $N(C=O)OR_5$

 NSO_2R_5 ,

NSO₂NR₅,

O,

S,

SO, or

SO₂NR₅,

and the other of X, Y, or Z is CH_2 or CR_3R_4 .

40. The compound of claim 39 as designated in formula IVF.

41. The compound of claim 39 as designated in formula IVG.

IVH

42. The compound of claim 38, wherein one of X, Y, or Z is $C=C-R_5$,

O=C,

10 NR₅,

 $N(C=O)R_5$,

 $N(C=O)OR_5$

NSO₂R₅,

NSO₂NR₅,

15 O,

S,

SO, or

SO₂NR₅,

and the others of X, Y, or Z is CH₂.

20

5

43. A compound of formula V

or a pharmaceutically acceptable salt thereof wherein:

5 A is O,

NH, or

S;

B is

10 $C(=O)R_1$,

 $C(=S)R_1$

heterocylco,

heteroaryl,

C(=O)-heterocyclo,

C(=N)-CN,or

C(=O)-heteteroaryl;

either D is N, E is C, and F is CH when "----" is a bond, or D is CH, E is N, and F is CH₂ when "----" is absent;

20

J, K, Q independently are CR_2 or N, with the proviso that when any one of J, K, or Q is N, then the other two are CR_2 ;

"---- "is absent; or is a bond; and

X, Y, Z independently are C=C-R₅,

25

O=C,

 CH_2 ,

CHR_{3.}

CHR₄,

CR₃R₄,

NR₅, $N(C=O)R_5$ $N(C=O)OR_5$, NSO₂R₅, 5 NSO₂NR₅, O, S, SO, or 10 SO₂; R₁ is H, (C_1-C_8) alkyl, (C₃-C₆)cycloalkyl, 15 O— $(C_1$ - C_4)alkyl, O—(C₃-C₆)cycloalkyl, S— $(C_1$ - $C_4)$ alkyl, S—(C₃-C₆)cycloalkyl, NH₂, 20 $NH(C_1-C_4)$ alkyl, $N((C_1-C_4)alkyl)_2$, or NH—(C₃-C₆)cycloalkyl, R₂ is H, 25 halo, (C₁-C₈)alkyl, (C₃-C₆)cycloalkyl, O— $(C_1$ - C_4)alkyl, O-(C₃-C₆)cycloalkyl, 30 S— (C_1-C_4) alkyl, S—(C₃-C₆)cycloalkyl, NH₂,

 $NH(C_1-C_4)$ alkyl,

 $N((C_1-C_4)alkyl)_2$, or NH—(C₃-C₆)cycloalkyl; 5 R₃ and R₄ independently are halo, (C_1-C_8) alkyl, (C₃-C₆)cycloalkyl, O— $(C_1$ - C_4)alkyl, O—(C₃-C₆)cycloalkyl, 10 S— $(C_1$ - $C_4)$ alkyl, S— $(C_3$ - $C_6)$ cycloalkyl, NH_2 , NH(C₁-C₄)alkyl, $N((C_1-C_4)alkyl)_2$ NH—(C₃-C₆)cycloalkyl; 15 aryl, $(CH_2)_n$ -aryl, heterocyclo, (CH₂)_n-heterocyclo, 20 heteroaryl, or (CH₂)_n-heteroaryl, wherein n is 0, 1, 2, or 3; R₅ is H, 25 (C_1-C_8) alkyl, (C₃-C₆)cycloalkyl, aryl, $(CH_2)_n$ -aryl, heterocyclo, 30 (CH₂)_n-heterocyclo,

heteroaryl, or

(CH₂)_n-heteroaryl,

wherein n is as defined above.

44. The compound of claim 43 as designated in formula VA.

VA

5

45. The compound of claim 43 as designated in formula VB.

10

46. The compound of claim 43 as designated in formula VC.

15 47. The compound of claim 43 as designated in formula VD

VD

wherein J_a is N or CR_6 , wherein R_6 is H or F.

5 48. The compound of claim 43 as designated in formula VE.

VE

49. The compound of claim 48, wherein two of X, Y, or Z is C=C-R₅,

10 O=C,

 NR_5 ,

 $N(C=O)R_5$

 $N(C=O)OR_5$

NSO₂R₅,

15 NSO₂NR₅,

O,

S,

SO, or

SO₂NR₅,

and the other of X, Y, or Z is CH_2 or CR_3R_4 .

50. The compound of claim 51 as designated in formula VF.

53. The compound of claim 51 as designated in formula VG.

5

54. The compound of claim 48, wherein one of X, Y, or Z is C=C-R₅,

O=C,

NR₅,

 $N(C=O)R_5$

10

15

 $N(C=O)OR_5$

NSO₂R₅,

NSO₂NR₅,

Ο,

S,

SO, or SO₂NR₅,

and the others of X, Y, or Z is CH₂.

/

55. A compound which is:

20 (S)-N-[2-Oxo-3-(6,7,8,9-tetrahydro-5H-benzocyclohepten-2-yl)-oxazolidin-5-ylmethyl]-acetamide;

(S)-N-[2-Oxo-3-(5-oxo-6,7,8,9-tetrahydro-5H-benzocyclohepten-2-yl)-oxazolidin-5-ylmethyl]-acetamide;

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(S)-N-[3-(6-Bromo-5-oxo-6,7,8,9-tetrahydro-5H-benzocyclohepten-2-yl)-2-oxo-oxazolidin-5-ylmethyl]-acetamide;

(S)-N-[3-(6-Dimethylaminomethylene-5-oxo-6,7,8,9-tetrahydro-5H-benzocyclohepten-2-yl)-2-oxo-oxazolidin-5-ylmethyl]-acetamide;

- (S)-N-[2-Oxo-3-(5-oxo-6-pyridin-4-ylmethylene-6,7,8,9-tetrahydro-5H-benzocyclohepten-2-yl)-oxazolidin-5-ylmethyl]-acetamide;
- (S)-N-[3-(6-Benzylidene-5-oxo-6,7,8,9-tetrahydro-5H-benzocyclohepten-5-oxo-oxazolidin-5-ylmethyl]-acetamide;
 - (S)-N-{3-[6-(4-Fluoro-benzylidene)-5-oxo-6,7,8,9-tetrahydro-5H-benzocyclohepten-2-yl]-2-oxo-oxazolidin-5-ylmethyl}-acetamide;
- (S)-N-[2-Oxo-3-(5-oxo-6-thiophen-3-ylmethylene-6,7,8,9-tetrahydro-5H-benzocyclohepten-2-yl)-oxazolidin-5-ylmethyl]-acetamide;
 - (S)-N-[3-(6-Furan-3-ylmethylene-5-oxo-6,7,8,9-tetrahydro-5H-benzocyclohepten-2-yl)-2-oxo-oxazolidin-5-ylmethyl]-acetamide;

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(S)-N-[2-Oxo-3-(6-oxo-6,7,8,9-tetrahydro-5H-benzocyclohepten-2-yl)-oxazolidin-5-ylmethyl]-acetamide;

- (S)-N-[2-Oxo-3-(7-oxo-6,7,8,9-tetrahydro-5H-benzocyclohepten-2-yl)-20 oxazolidin-5-ylmethyl]-acetamide;
 - (S)-N-[2-Oxo-3-(8-oxo-6,7,8,9-tetrahydro-5H-benzocyclohepten-2-yl)-oxazolidin-5-ylmethyl]-acetamide;
- 25 (S)-N-[2-Oxo-3-(9-oxo-6,7,8,9-tetrahydro-5H-benzocyclohepten-2-yl)-oxazolidin-5-ylmethyl]-acetamide;
- (S)-N-[3-(8,9-Dihydro-7H-benzocyclohepten-2-yl)-2-oxo-oxazolidin-5-ylmethyl]-acetamide;
 - (S)-N-[3-(8,9-Dihydro-5H-benzocyclohepten-2-yl)-2-oxo-oxazolidin-5-ylmethyl]-acetamide;
- (S)-N-[3-(6,9-Dihydro-5H-benzocyclohepten-2-yl)-2-oxo-oxazolidin-5-ylmethyl]-acetamide;
 - (S)-N-[3-(6,7-Dihydro-5H-benzocyclohepten-2-yl)-2-oxo-oxazolidin-5-ylmethyl]-acetamide;
- 40 (S)-N-[2-Oxo-3-(2,3,4,5-tetrahydro-1H-benzo[b]azepin-7-yl)-oxazolidin-5-ylmethyl]-acetamide;
 - (S)-N-[2-Oxo-3-(2,3,4,5-tetrahydro-1H-benzo[c]azepin-7-yl)-oxazolidin-5-ylmethyl]-acetamide;
 - (S)-N-[2-Oxo-3-(2,3,4,5-tetrahydro-1H-benzo[d]azepin-7-yl)-oxazolidin-5-ylmethyl]-acetamide;

- (S)-N-[2-Oxo-3-(2,3,4,5-tetrahydro-1H-benzo[c]azepin-8-yl)-oxazolidin-5-ylmethyl]-acetamide;
- 5 (S)-N-[2-Oxo-3-(2,3,4,5-tetrahydro-1H-benzo[b]azepin-8-yl)-oxazolidin-5-ylmethyl]-acetamide;
 - (S)-N-[2-Oxo-3-(2,3,4,5-tetrahydro-benzo[b]oxepin-7-yl)-oxazolidin-5-ylmethyl]-acetamide;
 - (S)-N-[2-Oxo-3-(1,3,4,5-tetrahydro-benzo[c]oxepin-7-yl)-oxazolidin-5-ylmethyl]-acetamide;
- (S)-N-[2-Oxo-3-(5,6,8,9-tetrahydro-7-oxa-benzocyclohepten-2-yl)oxazolidin-5-ylmethyl]-acetamide;

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- (S)-N-[2-Oxo-3-(1,3,4,5-tetrahydro-benzo[c]oxepin-8-yl)-oxazolidin-5-ylmethyl]-acetamide;
- 20 (S)-N-[2-Oxo-3-(2,3,4,5-tetrahydro-benzo[b]oxepin-8-yl)-oxazolidin-5-ylmethyl]-acetamide;
 - (S)-N-[2-Oxo-3-(6,7,8,9-tetrahydro-5-oxa-7-aza-benzocyclohepten-2-yl)-oxazolidin-5-ylmethyl]-acetamide;
 - (S)-N-[2-Oxo-3-(2,3,4,5-tetrahydro-benzo[b]thiepin-7-yl)-oxazolidin-5-ylmethyl]-acetamide;
- (S)-N-[2-Oxo-3-(1,3,4,5-tetrahydro-benzo[c]thiepin-7-yl)-oxazolidin-5-ylmethyl]-acetamide;
 - (S)-N-[2-Oxo-3-(1,2,4,5-tetrahydro-benzo[d]thiepin-7-yl)-oxazolidin-5-ylmethyl]-acetamide;
- 35 (S)-N-[2-Oxo-3-(1,3,4,5-tetrahydro-benzo[c]thiepin-8-yl)-oxazolidin-5-ylmethyl]-acetamide;
 - (S)-N-[2-Oxo-3-(2,3,4,5-tetrahydro-benzo[b]thiepin-8-yl)-oxazolidin-5-ylmethyl]-acetamide;
 - (S)-N-[2-Oxo-3-(5-oxo-2,3,4,5-tetrahydro-benzo[b]oxepin-8-yl)-oxazolidin-5-ylmethyl]-acetamide;
- (S)-N-[2-Oxo-3-(5-oxo-2,3,4,5-tetrahydro-benzo[b]thiepin-8-yl)oxazolidin-5-ylmethyl]-acetamide;

- (S)-N-[2-Oxo-3-(1,1,5-trioxo-2,3,4,5-tetrahydro-1H-1I6-benzo[b]thiepin-8-yl)-oxazolidin-5-ylmethyl]-acetamide;
- (S)-N-[2-Oxo-3-(5-oxo-2,3,4,5-tetrahydro-benzo[b]oxepin-7-yl)-5 oxazolidin-5-ylmethyl]-acetamide;
 - (S)-N-[3-(6,6-Difluoro-5-oxo-6,7,8,9-tetrahydro-5H-benzocyclohepten-2-yl)-2-oxo-oxazolidin-5-ylmethyl]-acetamide;
- 10 (S)-N-[3-(6-Benzylidene-5-oxo-6,7,8,9-tetrahydro-5H-benzocyclohepten-2-yl)-2-oxo-oxazolidin-5-ylmethyl]-acetamide;
 - (S)-N-[3-(2-Methyl-2,3,4,5-tetrahydro-1H-benzo[c]azepin-7-yl)-2-oxo-oxazolidin-5-ylmethyl]-acetamide; or
 - (S)-N-[3-(3-Methyl-2,3,4,5-tetrahydro-1H-benzo[d]azepin-7-yl)-2-oxo-oxazolidin-5-ylmethyl]-acetamide.
- 56. A pharmaceutical formulation comprising a compound of claim 1 admixedwith a pharmaceutically acceptable diluent, carrier, or excipient.
 - 57. A method of treating a bacterial infection in a mammal, comprising administering to a mammal in need thereof an effective amount of a compound of claim 1.

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